Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2021 – 06:36 PM EST

PDB ID	:	3K4H
Title	:	CRYSTAL STRUCTURE OF putative transcriptional regulator LacI from Bacillus cereus subsp. cytotoxis NVH 391-98
Authors	:	Malashkevich, V.N.; Toro, R.; Morano, C.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-10-05
Resolution	:	2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ⓘ) were used in the production of this report:

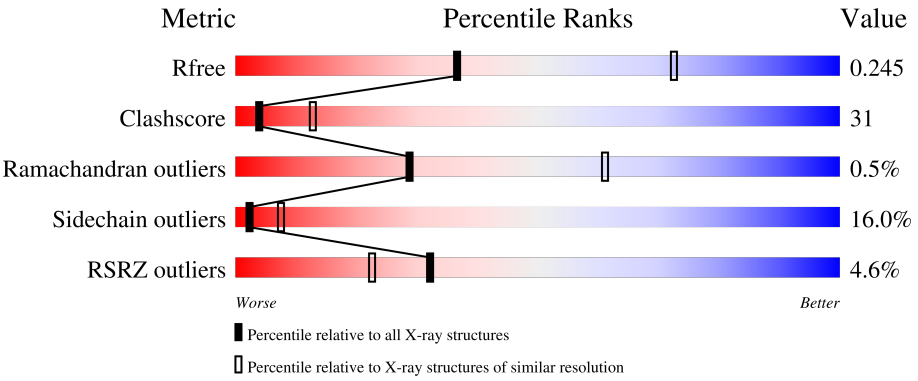
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	292	67% 26%
1	B	292	8% 42% 40% 12% 5%
2	C	2	50% 50%
2	D	2	50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	D	1	-	-	-	X
2	GLC	D	2	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4417 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called putative transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	282	Total	C	N	O	S	Se	0	0	0
			2188	1398	364	418	2	6			
1	B	278	Total	C	N	O	S	Se	0	0	0
			2159	1381	358	413	1	6			

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is water.

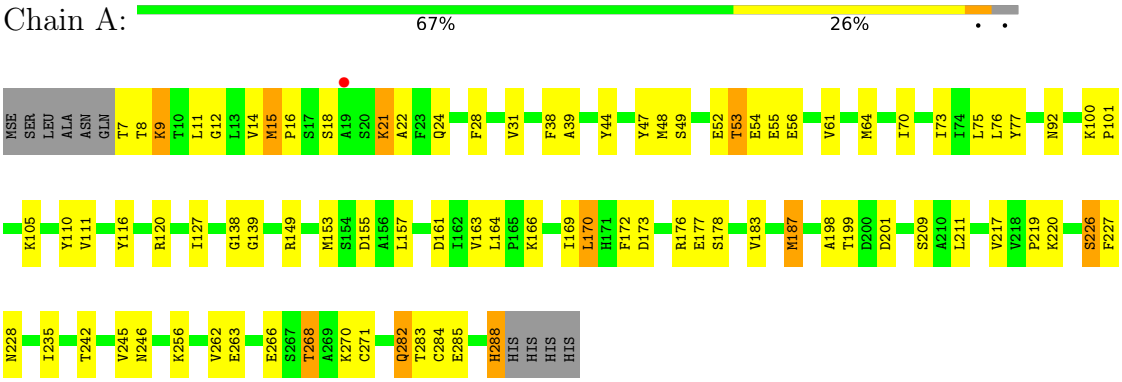
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	9	Total	O	0	0
			9	9		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

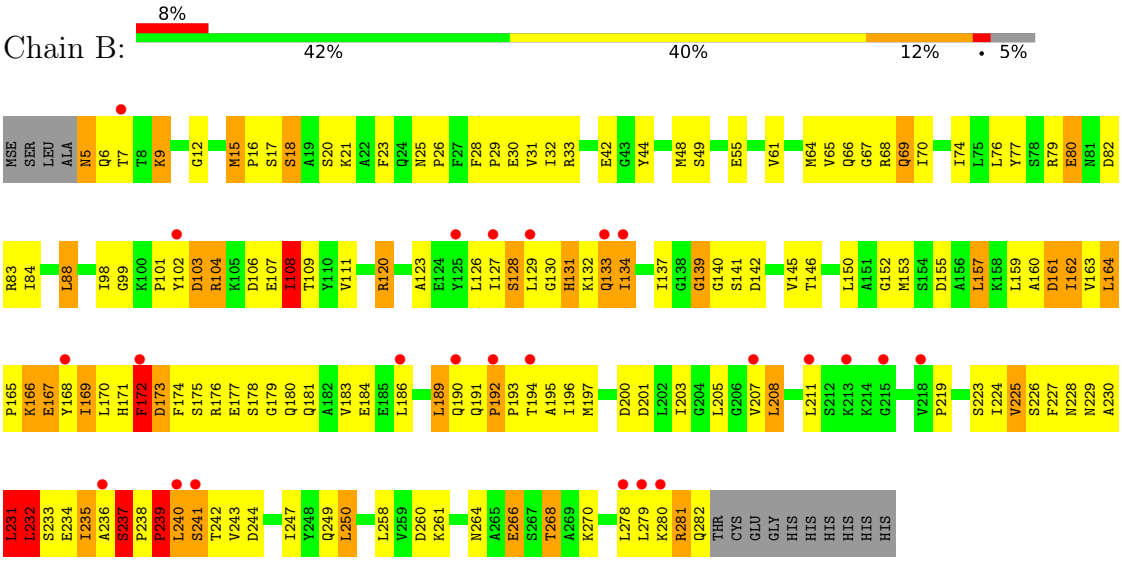
• Molecule 1: putative transcriptional regulator

Chain A:



• Molecule 1: putative transcriptional regulator

Chain B:



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C:

• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	98.61Å 98.61Å 186.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 39.88 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.80) 99.4 (39.88-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.98 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.203 , 0.244 0.210 , 0.245	Depositor DCC
R_{free} test set	1193 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	76.0	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 59.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4417	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GLC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	1.07	5/2223 (0.2%)	0.88	3/3000 (0.1%)
1	B	1.18	3/2192 (0.1%)	1.01	17/2958 (0.6%)
All	All	1.13	8/4415 (0.2%)	0.95	20/5958 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	240	LEU	CA-C	-17.10	1.08	1.52
1	B	240	LEU	C-O	-14.51	0.95	1.23
1	A	284	CYS	CB-SG	-8.39	1.68	1.82
1	B	241	SER	N-CA	-5.27	1.35	1.46
1	A	262	VAL	CB-CG1	-5.26	1.41	1.52
1	A	263	GLU	CG-CD	-5.25	1.44	1.51
1	A	111	VAL	CB-CG1	-5.11	1.42	1.52
1	A	111	VAL	CB-CG2	-5.08	1.42	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	240	LEU	N-CA-C	8.50	133.95	111.00
1	B	240	LEU	CA-C-O	-7.42	104.53	120.10
1	B	9	LYS	N-CA-C	7.11	130.19	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	240	LEU	O-C-N	6.85	133.66	122.70
1	B	169	ILE	CB-CA-C	-6.45	98.71	111.60
1	B	139	GLY	N-CA-C	-6.32	97.31	113.10
1	A	266	GLU	N-CA-C	-6.13	94.44	111.00
1	B	173	ASP	N-CA-C	6.08	127.42	111.00
1	B	240	LEU	CB-CG-CD1	-5.91	100.95	111.00
1	B	108	ILE	CB-CA-C	-5.87	99.86	111.60
1	B	172	PHE	N-CA-C	5.85	126.80	111.00
1	B	231	LEU	CA-CB-CG	-5.75	102.08	115.30
1	B	240	LEU	CB-CG-CD2	-5.69	101.32	111.00
1	B	232	LEU	N-CA-C	-5.62	95.84	111.00
1	A	170	LEU	CB-CG-CD2	-5.58	101.52	111.00
1	B	239	PRO	C-N-CA	5.49	135.42	121.70
1	B	250	LEU	CA-CB-CG	5.49	127.92	115.30
1	A	282	GLN	N-CA-C	5.33	125.40	111.00
1	B	240	LEU	N-CA-CB	-5.28	99.84	110.40
1	B	140	GLY	C-N-CA	5.14	134.55	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	231	LEU	Peptide
1	B	237	SER	Mainchain
1	B	239	PRO	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2188	0	2202	61	0
1	B	2159	0	2181	212	0
2	C	23	0	21	1	0
2	D	23	0	21	5	0
3	A	15	0	0	2	0
3	B	9	0	0	0	0
All	All	4417	0	4425	269	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

All (269) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:SER:HB3	1:B:238:PRO:CD	1.73	1.19
1:B:232:LEU:N	1:B:232:LEU:HD23	1.54	1.13
1:B:237:SER:CB	1:B:238:PRO:HD2	1.79	1.12
1:B:167:GLU:OE1	1:B:167:GLU:HA	1.53	1.08
1:B:12:GLY:HA3	1:B:64:MSE:HE1	1.29	1.07
1:B:237:SER:HB3	1:B:238:PRO:HD2	1.26	1.06
1:B:104:ARG:HG3	1:B:104:ARG:HH11	1.19	1.06
1:B:230:ALA:O	1:B:231:LEU:HD23	1.56	1.05
1:B:165:PRO:O	1:B:166:LYS:HB2	1.53	1.03
1:B:169:ILE:HG22	1:B:169:ILE:O	1.57	1.01
1:B:131:HIS:NE2	1:B:194:THR:HG22	1.75	1.00
1:B:131:HIS:HE2	1:B:194:THR:HG22	1.24	0.99
1:B:131:HIS:NE2	1:B:194:THR:O	1.96	0.98
1:B:231:LEU:O	1:B:231:LEU:HD12	1.60	0.98
1:B:230:ALA:CB	1:B:232:LEU:HG	1.93	0.98
1:A:176:ARG:HH11	1:A:176:ARG:HG3	1.29	0.98
1:A:227:PHE:O	1:A:228:ASN:HB2	1.64	0.98
1:B:232:LEU:H	1:B:232:LEU:HD23	1.29	0.98
1:B:69:GLN:HA	1:B:69:GLN:HE21	1.27	0.97
1:B:15:MSE:HE1	1:B:48:MSE:SE	2.15	0.97
1:A:12:GLY:HA3	1:A:64:MSE:HE1	1.48	0.96
1:B:104:ARG:CG	1:B:104:ARG:HH11	1.78	0.95
1:B:164:LEU:CD2	1:B:169:ILE:HD11	1.97	0.94
1:B:64:MSE:HE3	1:B:70:ILE:HD13	1.52	0.91
1:B:162:ILE:CD1	1:B:162:ILE:N	2.32	0.91
1:B:133:GLN:NE2	1:B:168:TYR:OH	2.03	0.91
1:B:230:ALA:HB1	1:B:232:LEU:HG	1.51	0.91
1:B:172:PHE:N	1:B:172:PHE:CD2	2.30	0.89
1:B:12:GLY:HA3	1:B:64:MSE:CE	2.01	0.89
1:A:139:GLY:HA2	1:A:172:PHE:O	1.73	0.89
1:B:194:THR:O	1:B:194:THR:HG22	1.73	0.88
1:B:172:PHE:H	1:B:172:PHE:HD2	1.17	0.87
1:B:164:LEU:HD23	1:B:169:ILE:HD11	1.59	0.84
1:B:225:VAL:HA	1:B:241:SER:HB2	1.57	0.84
1:B:127:ILE:O	1:B:130:GLY:N	2.11	0.84
1:B:139:GLY:HA3	1:B:174:PHE:CE1	2.13	0.84
1:B:127:ILE:O	1:B:130:GLY:HA2	1.78	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:ILE:O	1:B:130:GLY:CA	2.26	0.83
1:B:162:ILE:HD13	1:B:162:ILE:H	1.44	0.82
1:B:237:SER:CB	1:B:238:PRO:CD	2.41	0.82
1:A:282:GLN:NE2	3:A:522:HOH:O	2.13	0.81
1:B:232:LEU:N	1:B:232:LEU:CD2	2.30	0.81
1:B:193:PRO:O	1:B:194:THR:HB	1.79	0.80
1:B:165:PRO:O	1:B:166:LYS:CB	2.29	0.80
1:B:230:ALA:O	1:B:231:LEU:CD2	2.30	0.80
1:B:162:ILE:N	1:B:162:ILE:HD13	1.97	0.80
1:B:99:GLY:CA	2:D:2:GLC:H3	2.12	0.79
1:B:225:VAL:O	1:B:225:VAL:CG1	2.30	0.79
1:B:17:SER:HB3	1:B:20:SER:HB3	1.64	0.79
1:B:231:LEU:O	1:B:231:LEU:CD1	2.30	0.79
1:B:165:PRO:HG3	1:B:168:TYR:CD1	2.17	0.79
1:B:165:PRO:HB2	1:B:167:GLU:H	1.49	0.78
1:B:231:LEU:HA	1:B:234:GLU:OE2	1.84	0.78
1:B:157:LEU:HD23	1:B:164:LEU:HD13	1.66	0.77
1:B:170:LEU:O	1:B:171:HIS:ND1	2.17	0.77
1:B:67:GLY:O	1:B:68:ARG:HB2	1.83	0.77
1:B:99:GLY:HA3	2:D:2:GLC:H3	1.66	0.77
1:A:176:ARG:NH1	1:A:176:ARG:HG3	1.99	0.77
1:B:167:GLU:OE1	1:B:167:GLU:CA	2.30	0.76
1:B:9:LYS:NZ	1:B:42:GLU:O	2.19	0.76
1:B:237:SER:HB3	1:B:238:PRO:HD3	1.67	0.75
1:B:230:ALA:HB3	1:B:232:LEU:HG	1.65	0.75
1:B:104:ARG:HD3	1:B:107:GLU:OE1	1.87	0.74
1:B:225:VAL:HG23	1:B:241:SER:HB3	1.69	0.74
1:B:164:LEU:HD21	1:B:169:ILE:HD11	1.67	0.74
1:B:170:LEU:HD11	1:B:186:LEU:HD13	1.68	0.74
1:B:131:HIS:CD2	1:B:194:THR:O	2.41	0.74
1:B:15:MSE:HG3	1:B:16:PRO:HD2	1.69	0.73
1:B:131:HIS:CD2	1:B:194:THR:HG22	2.22	0.72
1:B:104:ARG:HG3	1:B:104:ARG:NH1	1.95	0.72
1:B:225:VAL:O	1:B:225:VAL:HG12	1.88	0.72
1:B:237:SER:HB2	1:B:238:PRO:HD2	1.71	0.72
1:B:189:LEU:HG	1:B:190:GLN:H	1.54	0.72
1:B:162:ILE:HD12	1:B:162:ILE:N	2.05	0.72
1:B:165:PRO:HG3	1:B:168:TYR:CG	2.25	0.71
1:B:224:ILE:O	1:B:241:SER:HB2	1.89	0.71
1:B:208:LEU:HD13	1:B:237:SER:HB2	1.72	0.71
1:A:201:ASP:OD1	1:A:226:SER:OG	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:GLU:C	1:A:54:GLU:OE2	2.29	0.70
1:B:207:VAL:O	1:B:211:LEU:HG	1.93	0.69
1:B:230:ALA:O	1:B:231:LEU:CB	2.41	0.69
1:B:281:ARG:O	1:B:282:GLN:C	2.30	0.68
1:A:173:ASP:OD1	1:A:178:SER:OG	2.10	0.67
1:A:12:GLY:HA3	1:A:64:MSE:CE	2.24	0.67
1:B:165:PRO:HG2	1:B:168:TYR:HB2	1.75	0.67
1:B:157:LEU:CD2	1:B:164:LEU:HD13	2.25	0.67
1:A:127:ILE:HD11	1:A:157:LEU:HD23	1.77	0.66
1:B:266:GLU:OE2	1:B:268:THR:N	2.28	0.66
1:B:164:LEU:CD2	1:B:169:ILE:CD1	2.74	0.66
1:B:227:PHE:O	1:B:228:ASN:CB	2.44	0.66
1:A:227:PHE:O	1:A:228:ASN:CB	2.41	0.65
1:B:196:ILE:CG2	1:B:197:MSE:N	2.59	0.65
1:B:64:MSE:CE	1:B:70:ILE:HD13	2.25	0.65
1:B:157:LEU:HD23	1:B:164:LEU:CD1	2.25	0.65
1:B:108:ILE:CG2	1:B:109:THR:N	2.57	0.65
1:A:173:ASP:C	1:A:173:ASP:OD1	2.32	0.64
1:A:53:THR:HG23	1:A:56:GLU:CD	2.17	0.64
1:B:123:ALA:O	1:B:127:ILE:HG12	1.98	0.64
1:B:133:GLN:HB2	1:B:193:PRO:HA	1.80	0.64
1:B:229:ASN:CA	1:B:242:THR:HG21	2.27	0.64
1:B:194:THR:CG2	1:B:194:THR:O	2.45	0.63
1:B:231:LEU:HG	1:B:232:LEU:HD23	1.80	0.63
1:B:104:ARG:NH1	1:B:104:ARG:CG	2.50	0.63
1:B:230:ALA:O	1:B:231:LEU:HB3	1.98	0.63
1:B:240:LEU:HD22	1:B:241:SER:H	1.63	0.63
1:B:160:ALA:O	1:B:161:ASP:CB	2.47	0.62
1:B:164:LEU:HD23	1:B:169:ILE:CD1	2.28	0.62
1:B:28:PHE:HA	1:B:31:VAL:HG12	1.81	0.61
1:B:16:PRO:HA	1:B:77:TYR:CE2	2.35	0.61
1:B:230:ALA:O	1:B:231:LEU:CG	2.49	0.61
1:B:157:LEU:CD2	1:B:164:LEU:CD1	2.79	0.60
1:A:64:MSE:CE	1:A:70:ILE:HG21	2.32	0.60
1:B:165:PRO:CG	1:B:168:TYR:CG	2.85	0.59
1:A:9:LYS:HB3	1:A:44:TYR:CE2	2.37	0.59
1:B:16:PRO:HA	1:B:77:TYR:HE2	1.67	0.59
1:B:229:ASN:HA	1:B:242:THR:HG21	1.85	0.58
1:B:227:PHE:O	1:B:228:ASN:HB2	2.03	0.58
1:A:288:HIS:ND1	1:A:288:HIS:N	2.51	0.58
1:B:162:ILE:HG22	1:B:163:VAL:N	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LEU:O	1:B:160:ALA:HB3	2.04	0.58
1:B:139:GLY:HA3	1:B:174:PHE:CZ	2.39	0.58
1:B:193:PRO:O	1:B:194:THR:CB	2.45	0.58
1:B:131:HIS:CD2	1:B:195:ALA:HB2	2.39	0.57
1:B:231:LEU:N	1:B:233:SER:H	2.01	0.57
1:B:111:VAL:HG12	1:B:250:LEU:HD22	1.86	0.57
1:B:226:SER:O	1:B:242:THR:HA	2.04	0.57
1:B:69:GLN:CA	1:B:69:GLN:HE21	1.99	0.57
1:A:268:THR:HG21	1:B:249:GLN:NE2	2.19	0.57
1:B:137:ILE:HG23	1:B:172:PHE:HE2	1.70	0.57
1:B:160:ALA:O	1:B:161:ASP:HB2	2.03	0.57
1:B:127:ILE:C	1:B:130:GLY:H	2.07	0.57
1:B:225:VAL:HG23	1:B:241:SER:CB	2.34	0.57
1:B:232:LEU:H	1:B:232:LEU:CD2	1.94	0.56
1:A:256:LYS:NZ	1:B:266:GLU:HG3	2.21	0.56
1:B:103:ASP:OD1	1:B:103:ASP:N	2.30	0.55
1:B:165:PRO:HG2	1:B:168:TYR:CB	2.36	0.55
1:A:169:ILE:O	1:A:170:LEU:HD23	2.05	0.55
1:A:18:SER:CB	1:A:52:GLU:HG2	2.36	0.55
1:A:226:SER:OG	1:A:227:PHE:N	2.37	0.55
1:B:7:THR:HG23	1:B:7:THR:O	2.05	0.55
1:B:16:PRO:CA	1:B:77:TYR:HE2	2.20	0.54
1:B:99:GLY:HA2	2:D:2:GLC:H3	1.89	0.54
1:B:61:VAL:O	1:B:65:VAL:HG23	2.08	0.54
1:B:133:GLN:HB3	1:B:168:TYR:HE2	1.72	0.53
1:B:228:ASN:O	1:B:229:ASN:C	2.43	0.53
1:B:155:ASP:O	1:B:159:LEU:HG	2.08	0.53
1:B:164:LEU:HD21	1:B:169:ILE:CD1	2.36	0.53
1:A:18:SER:HB2	1:A:52:GLU:HG2	1.90	0.53
1:A:177:GLU:CD	1:A:177:GLU:H	2.10	0.53
1:A:15:MSE:HG2	1:A:22:ALA:HB2	1.92	0.52
1:B:111:VAL:HG12	1:B:250:LEU:CD2	2.40	0.52
1:B:169:ILE:O	1:B:170:LEU:HD23	2.09	0.52
1:B:18:SER:HA	1:B:23:PHE:CD1	2.44	0.52
1:A:64:MSE:HE1	1:A:73:ILE:HG12	1.92	0.51
1:A:16:PRO:HG3	1:A:77:TYR:HB3	1.91	0.51
1:B:145:VAL:HG13	1:B:146:THR:N	2.25	0.51
1:A:14:VAL:HB	1:A:75:LEU:HD23	1.91	0.51
1:B:133:GLN:C	1:B:134:ILE:HG12	2.27	0.50
1:A:173:ASP:O	1:A:173:ASP:OD1	2.30	0.50
1:B:157:LEU:HG	1:B:162:ILE:HB	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:VAL:O	1:A:64:MSE:HB2	2.12	0.50
1:B:153:MSE:CE	1:B:197:MSE:HB2	2.41	0.50
1:B:69:GLN:HA	1:B:69:GLN:NE2	2.11	0.50
1:A:226:SER:O	1:A:242:THR:HA	2.12	0.49
1:B:172:PHE:O	1:B:173:ASP:OD2	2.30	0.49
1:B:108:ILE:HG23	1:B:109:THR:N	2.26	0.49
1:B:229:ASN:O	1:B:229:ASN:OD1	2.29	0.49
1:B:153:MSE:HE3	1:B:197:MSE:HB2	1.92	0.49
1:B:186:LEU:HD23	1:B:196:ILE:HD11	1.95	0.49
1:B:197:MSE:HG3	1:B:225:VAL:CG1	2.43	0.49
1:A:138:GLY:HA2	1:A:199:THR:OG1	2.12	0.49
1:B:266:GLU:OE2	1:B:268:THR:OG1	2.31	0.49
1:B:169:ILE:C	1:B:170:LEU:HG	2.32	0.49
1:B:98:ILE:HG22	2:D:2:GLC:O3	2.12	0.49
1:B:278:LEU:HD12	1:B:278:LEU:C	2.34	0.48
1:B:229:ASN:HD22	1:B:244:ASP:HB2	1.78	0.48
1:B:231:LEU:C	1:B:233:SER:N	2.62	0.48
1:A:53:THR:HG23	1:A:56:GLU:OE1	2.13	0.48
1:B:108:ILE:HG22	1:B:109:THR:N	2.23	0.47
1:B:28:PHE:O	1:B:32:ILE:HG12	2.14	0.47
1:A:170:LEU:HA	1:A:170:LEU:HD23	1.53	0.47
1:B:231:LEU:H	1:B:233:SER:H	1.62	0.47
1:B:200:ASP:HB3	1:B:203:ILE:HG13	1.96	0.47
1:B:238:PRO:HG2	1:B:238:PRO:O	2.15	0.47
1:B:230:ALA:HB1	1:B:232:LEU:CG	2.33	0.47
1:B:30:GLU:OE2	1:B:33:ARG:NH1	2.48	0.47
1:B:5:ASN:HB3	1:B:6:GLN:H	1.61	0.46
1:B:196:ILE:HG23	1:B:197:MSE:N	2.29	0.46
1:B:69:GLN:CA	1:B:69:GLN:NE2	2.70	0.46
1:A:15:MSE:SE	1:A:48:MSE:HE1	2.66	0.46
1:A:220:LYS:HA	1:A:285:GLU:OE2	2.15	0.46
1:B:79:ARG:O	1:B:82:ASP:HB2	2.15	0.46
1:A:15:MSE:SE	1:A:48:MSE:CE	3.14	0.46
1:B:175:SER:OG	1:B:178:SER:HB3	2.16	0.46
1:B:196:ILE:HG22	1:B:197:MSE:N	2.28	0.46
1:B:104:ARG:O	1:B:107:GLU:N	2.48	0.46
1:B:280:LYS:HG3	1:B:280:LYS:O	2.15	0.46
1:B:80:GLU:OE2	1:B:102:TYR:HB2	2.16	0.46
1:B:172:PHE:O	1:B:173:ASP:CG	2.54	0.45
1:B:208:LEU:CD1	1:B:237:SER:HB2	2.42	0.45
1:A:38:PHE:CE2	1:A:256:LYS:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:LYS:O	1:A:21:LYS:HD2	2.16	0.45
1:A:15:MSE:HE2	1:A:28:PHE:CD2	2.52	0.45
1:B:238:PRO:HA	1:B:239:PRO:HD2	1.54	0.45
1:B:233:SER:OG	1:B:234:GLU:N	2.48	0.45
1:A:64:MSE:HE3	1:A:70:ILE:HG13	1.98	0.45
1:B:141:SER:HB3	1:B:171:HIS:CD2	2.53	0.44
1:B:162:ILE:CG2	1:B:163:VAL:N	2.81	0.44
1:B:242:THR:CG2	1:B:243:VAL:N	2.80	0.44
1:B:223:SER:OG	1:B:282:GLN:NE2	2.49	0.44
1:B:238:PRO:CG	1:B:238:PRO:O	2.65	0.44
1:A:219:PRO:HB3	1:A:283:THR:C	2.37	0.44
1:A:64:MSE:HE3	1:A:70:ILE:HG21	2.00	0.44
1:B:128:SER:C	1:B:130:GLY:N	2.68	0.44
1:B:224:ILE:O	1:B:241:SER:CB	2.61	0.44
1:A:18:SER:HB3	1:A:52:GLU:HG2	2.00	0.44
1:B:179:GLY:O	1:B:183:VAL:HG23	2.18	0.44
1:B:74:ILE:HG13	1:B:258:LEU:HD22	1.99	0.44
1:B:205:LEU:CD2	1:B:236:ALA:HB2	2.48	0.44
1:A:24:GLN:HB2	1:A:24:GLN:HE21	1.64	0.43
1:B:260:ASP:OD1	1:B:264:ASN:ND2	2.44	0.43
1:A:155:ASP:OD2	1:B:120:ARG:NH2	2.45	0.43
1:A:217:VAL:CG1	1:A:220:LYS:HG2	2.49	0.43
1:B:165:PRO:HG2	1:B:168:TYR:H	1.84	0.43
1:B:186:LEU:CD2	1:B:196:ILE:HD11	2.48	0.43
1:B:99:GLY:CA	2:D:2:GLC:C3	2.91	0.43
1:B:28:PHE:O	1:B:29:PRO:C	2.54	0.43
1:B:167:GLU:O	1:B:167:GLU:CG	2.67	0.43
1:A:15:MSE:HG2	1:A:22:ALA:CB	2.49	0.42
1:A:116:TYR:OH	1:B:120:ARG:NH2	2.52	0.42
1:B:16:PRO:CA	1:B:77:TYR:CE2	3.00	0.42
1:B:229:ASN:N	1:B:242:THR:HG21	2.34	0.42
1:B:25:ASN:HA	1:B:26:PRO:HD3	1.94	0.42
1:B:120:ARG:HG2	1:B:152:GLY:O	2.18	0.42
1:B:231:LEU:CA	1:B:233:SER:H	2.32	0.42
1:B:172:PHE:O	1:B:178:SER:HB2	2.19	0.42
1:B:134:ILE:HG23	1:B:134:ILE:HD13	1.74	0.42
1:B:228:ASN:OD1	1:B:247:ILE:HD11	2.20	0.42
1:A:198:ALA:O	1:A:226:SER:HA	2.20	0.42
1:A:101:PRO:HD3	1:A:110:TYR:CD2	2.55	0.42
1:A:55:GLU:HB2	3:A:527:HOH:O	2.20	0.42
1:B:26:PRO:HG2	1:B:232:LEU:HD11	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:HIS:CD2	1:B:194:THR:CG2	3.00	0.42
1:B:192:PRO:HA	1:B:193:PRO:HD3	1.75	0.41
1:B:235:ILE:O	1:B:235:ILE:HG23	2.20	0.41
1:B:238:PRO:HB2	1:B:282:GLN:O	2.20	0.41
1:B:195:ALA:O	1:B:196:ILE:HG13	2.20	0.41
1:B:197:MSE:HG3	1:B:225:VAL:HG11	2.03	0.41
1:B:172:PHE:O	1:B:178:SER:CB	2.68	0.41
1:B:235:ILE:HG22	1:B:236:ALA:N	2.33	0.41
1:A:11:LEU:HA	1:A:11:LEU:HD23	1.66	0.41
1:A:183:VAL:O	1:A:187:MSE:HB2	2.21	0.41
1:A:176:ARG:NH1	1:A:176:ARG:CG	2.74	0.41
1:A:38:PHE:O	1:A:39:ALA:C	2.58	0.41
1:A:47:TYR:CE2	1:A:64:MSE:HG2	2.56	0.41
1:B:9:LYS:O	1:B:44:TYR:HA	2.21	0.41
1:A:100:LYS:HA	1:A:101:PRO:HD3	1.94	0.41
1:B:165:PRO:CG	1:B:168:TYR:CD1	2.96	0.41
1:B:177:GLU:HG2	1:B:181:GLN:NE2	2.35	0.41
1:B:219:PRO:HB3	1:B:282:GLN:HE21	1.86	0.41
1:A:149:ARG:HH12	2:C:1:GLC:HO6	1.68	0.41
1:B:175:SER:OG	1:B:178:SER:CB	2.69	0.40
1:B:126:LEU:O	1:B:130:GLY:N	2.50	0.40
1:A:54:GLU:OE2	1:A:55:GLU:N	2.54	0.40
1:B:84:ILE:HG22	1:B:88:LEU:HD22	2.02	0.40
1:B:126:LEU:HD23	1:B:126:LEU:HA	1.88	0.40
1:A:64:MSE:CE	1:A:73:ILE:HG12	2.51	0.40
1:B:126:LEU:O	1:B:129:LEU:HB2	2.21	0.40
1:B:157:LEU:O	1:B:161:ASP:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	280/292 (96%)	277 (99%)	3 (1%)	0	100 100
1	B	276/292 (94%)	263 (95%)	10 (4%)	3 (1%)	14 41
All	All	556/584 (95%)	540 (97%)	13 (2%)	3 (0%)	29 61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	237	SER
1	B	101	PRO
1	B	192	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	239/241 (99%)	211 (88%)	28 (12%)	5 16
1	B	236/241 (98%)	188 (80%)	48 (20%)	1 4
All	All	475/482 (98%)	399 (84%)	76 (16%)	2 7

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	8	THR
1	A	9	LYS
1	A	15	MSE
1	A	21	LYS
1	A	31	VAL
1	A	49	SER
1	A	53	THR
1	A	76	LEU
1	A	92	ASN
1	A	105	LYS
1	A	120	ARG
1	A	153	MSE

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Mol	Chain	Res	Type
1	A	161	ASP
1	A	163	VAL
1	A	164	LEU
1	A	166	LYS
1	A	187	MSE
1	A	209	SER
1	A	211	LEU
1	A	226	SER
1	A	235	ILE
1	A	245	VAL
1	A	246	ASN
1	A	268	THR
1	A	270	LYS
1	A	271	CYS
1	A	288	HIS
1	B	5	ASN
1	B	15	MSE
1	B	18	SER
1	B	21	LYS
1	B	49	SER
1	B	55	GLU
1	B	66	GLN
1	B	69	GLN
1	B	76	LEU
1	B	80	GLU
1	B	83	ARG
1	B	88	LEU
1	B	103	ASP
1	B	104	ARG
1	B	106	ASP
1	B	108	ILE
1	B	120	ARG
1	B	128	SER
1	B	131	HIS
1	B	132	LYS
1	B	133	GLN
1	B	134	ILE
1	B	142	ASP
1	B	150	LEU
1	B	157	LEU
1	B	161	ASP
1	B	162	ILE

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Mol	Chain	Res	Type
1	B	164	LEU
1	B	166	LYS
1	B	167	GLU
1	B	172	PHE
1	B	176	ARG
1	B	180	GLN
1	B	184	GLU
1	B	189	LEU
1	B	191	GLN
1	B	201	ASP
1	B	208	LEU
1	B	225	VAL
1	B	231	LEU
1	B	232	LEU
1	B	235	ILE
1	B	261	LYS
1	B	266	GLU
1	B	268	THR
1	B	270	LYS
1	B	279	LEU
1	B	281	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	228	ASN
1	A	249	GLN
1	B	5	ASN
1	B	69	GLN
1	B	133	GLN
1	B	228	ASN
1	B	229	ASN
1	B	249	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

4 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	GLC	C	1	2	12,12,12	0.51	0	17,17,17	1.49	3 (17%)
2	GLC	C	2	2	11,11,12	0.69	0	15,15,17	1.72	5 (33%)
2	GLC	D	1	2	12,12,12	0.53	0	17,17,17	1.48	3 (17%)
2	GLC	D	2	2	11,11,12	0.87	0	15,15,17	2.09	6 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	GLC	C1-O5-C5	3.97	121.16	113.66
2	D	2	GLC	C1-O5-C5	3.67	117.16	112.19
2	D	2	GLC	C1-C2-C3	-3.32	105.59	109.67
2	C	2	GLC	O5-C5-C6	3.12	112.10	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	2	GLC	C2-C3-C4	3.08	116.23	110.89
2	C	2	GLC	O4-C4-C5	-2.98	101.89	109.30
2	C	1	GLC	C4-C3-C2	-2.83	105.89	110.82
2	D	2	GLC	O3-C3-C4	-2.74	104.02	110.35
2	C	1	GLC	O5-C1-C2	2.71	115.13	110.28
2	D	1	GLC	C6-C5-C4	-2.69	106.71	113.00
2	C	2	GLC	C1-C2-C3	-2.41	106.71	109.67
2	C	2	GLC	C1-O5-C5	2.29	115.30	112.19
2	D	1	GLC	O5-C5-C6	2.29	112.12	106.44
2	C	2	GLC	O3-C3-C4	2.22	115.47	110.35
2	D	1	GLC	C4-C3-C2	-2.09	107.17	110.82
2	D	2	GLC	O3-C3-C2	-2.08	106.00	109.99
2	D	2	GLC	O4-C4-C5	2.04	114.36	109.30

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	GLC	O5-C5-C6-O6
2	C	2	GLC	O5-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6
2	D	2	GLC	C4-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6

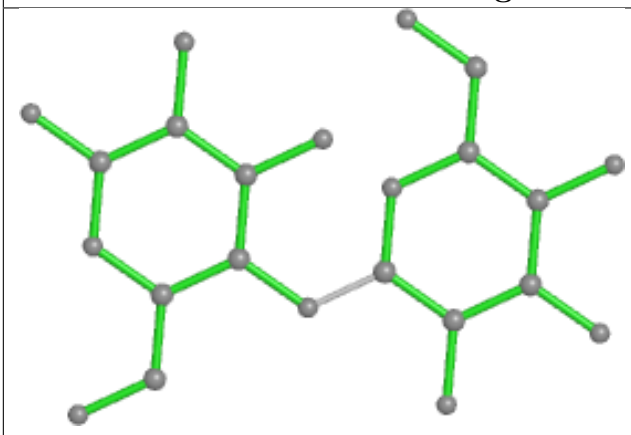
There are no ring outliers.

2 monomers are involved in 6 short contacts:

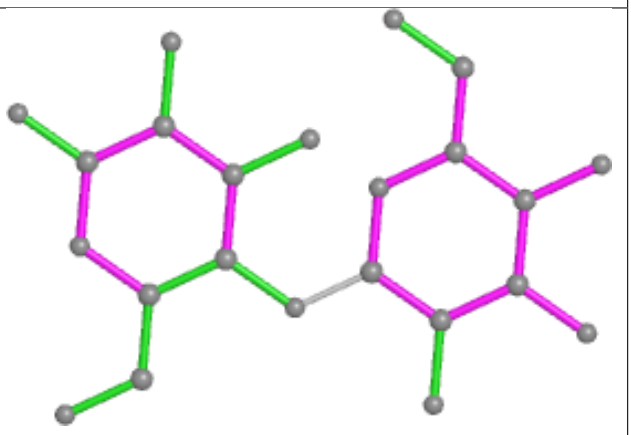
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	GLC	5	0
2	C	1	GLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

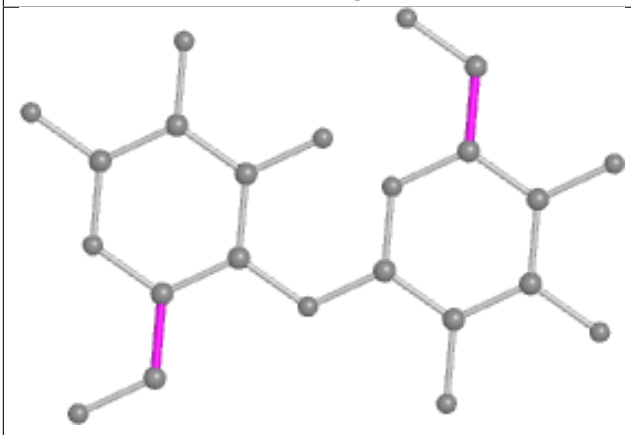
Oligosaccharide Chain C



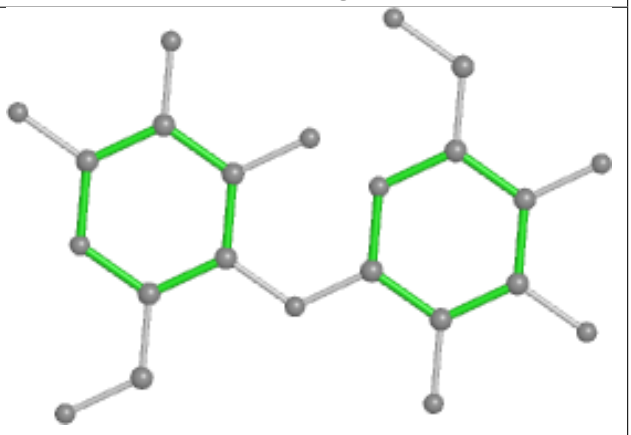
Bond lengths



Bond angles

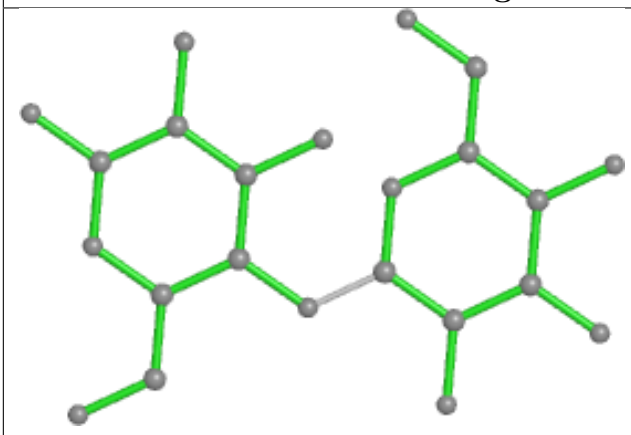


Torsions

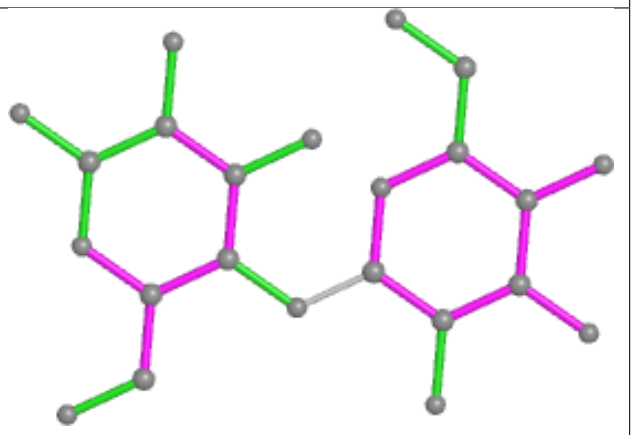


Rings

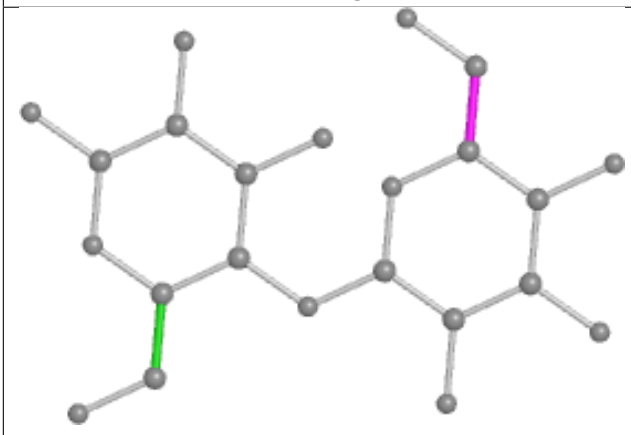
Oligosaccharide Chain D



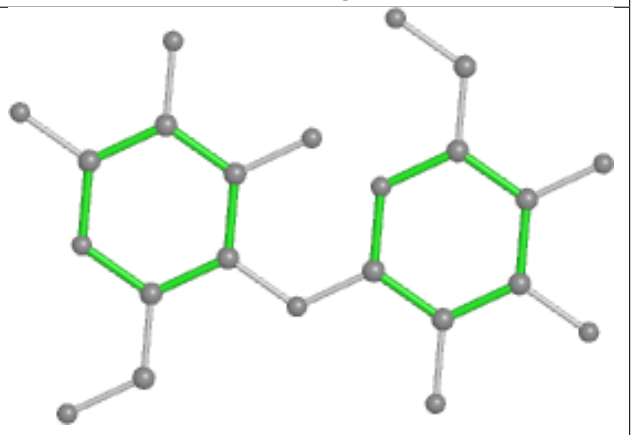
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	276/292 (94%)	-0.19	1 (0%) 92 91	36, 49, 73, 83	0
1	B	272/292 (93%)	0.36	24 (8%) 10 5	43, 64, 88, 94	0
All	All	548/584 (93%)	0.08	25 (4%) 32 22	36, 55, 85, 94	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	129	LEU	4.5
1	B	236	ALA	3.8
1	B	211	LEU	3.6
1	B	215	GLY	3.5
1	B	207	VAL	3.4
1	B	172	PHE	3.1
1	B	133	GLN	3.1
1	B	190	GLN	3.0
1	B	134	ILE	2.7
1	B	278	LEU	2.7
1	B	7	THR	2.5
1	B	280	LYS	2.5
1	B	102	TYR	2.5
1	B	127	ILE	2.5
1	B	186	LEU	2.4
1	B	192	PRO	2.3
1	B	213	LYS	2.3
1	B	241	SER	2.3
1	B	240	LEU	2.3
1	B	279	LEU	2.2
1	B	168	TYR	2.2
1	B	218	VAL	2.2
1	B	125	TYR	2.1
1	B	194	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	19	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

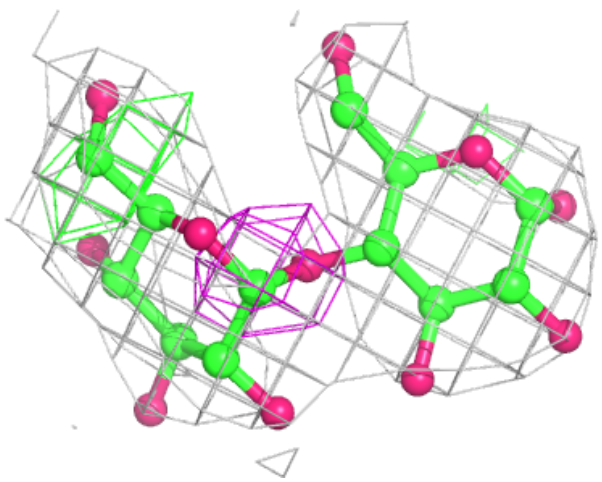
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

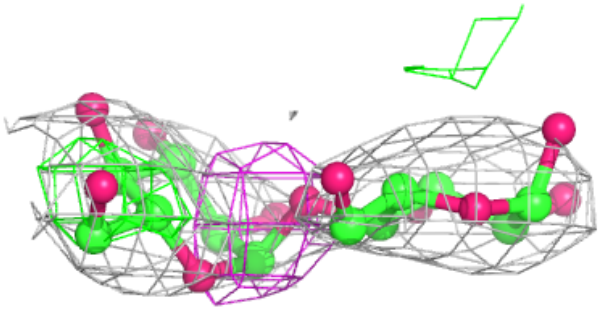
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GLC	D	2	11/12	0.59	0.60	46,47,52,53	11
2	GLC	D	1	12/12	0.68	0.72	45,45,52,52	12
2	GLC	C	2	11/12	0.81	0.32	62,64,67,67	11
2	GLC	C	1	12/12	0.86	0.29	65,65,67,67	12

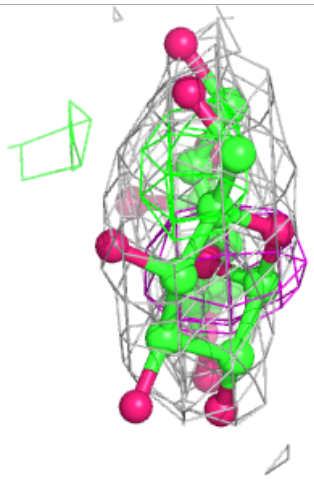
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

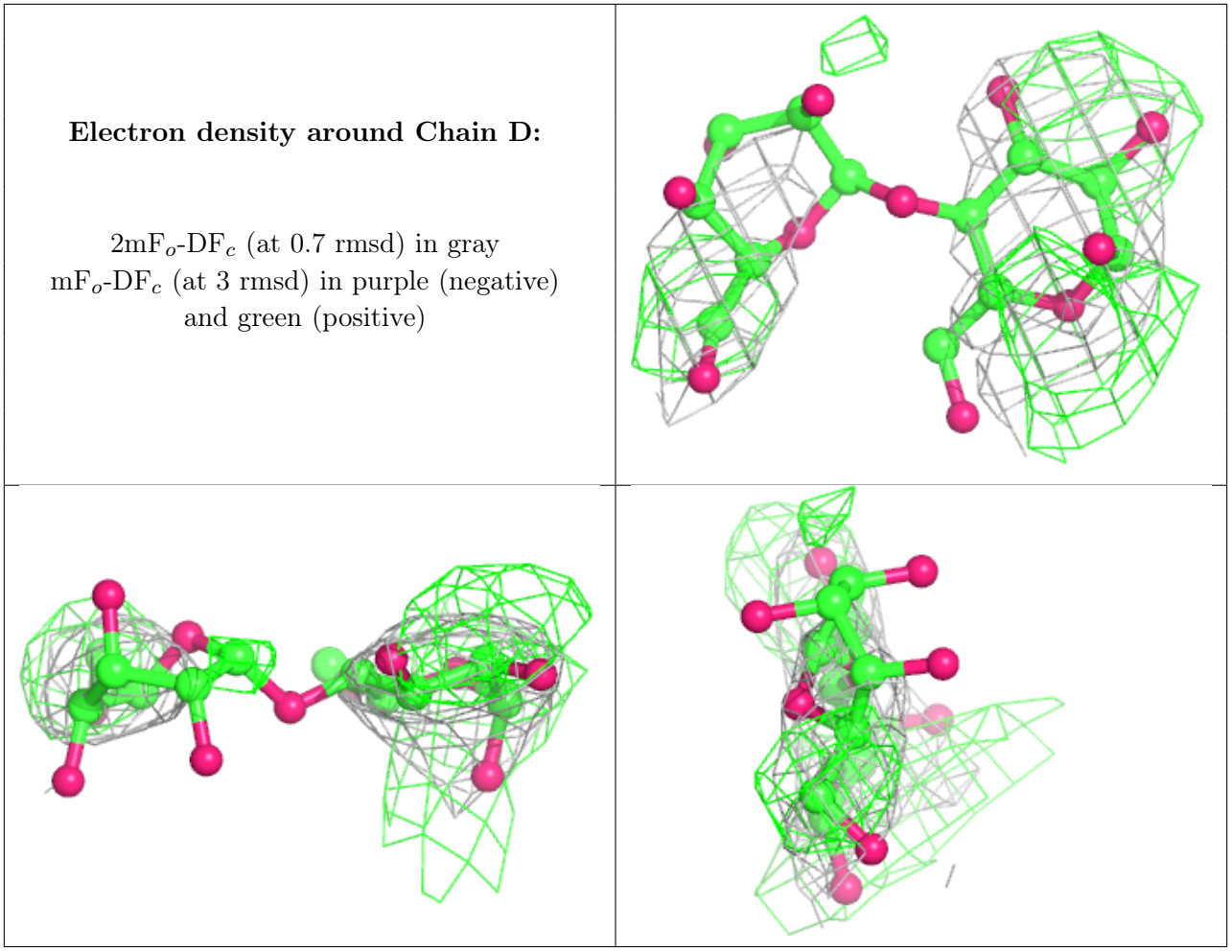






Electron density around Chain D:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.